



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:28 PM GMT

PDB ID : 1XF6  
Title : High resolution crystal structure of phycoerythrin 545 from the marine cryptophyte rhodomonas CS24  
Authors : Doust, A.B.; Marai, C.N.J.; Harrop, S.J.; Wilk, K.E.; Curmi, P.M.G.; Scholes, G.D.  
Deposited on : 2004-09-14  
Resolution : 1.10 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

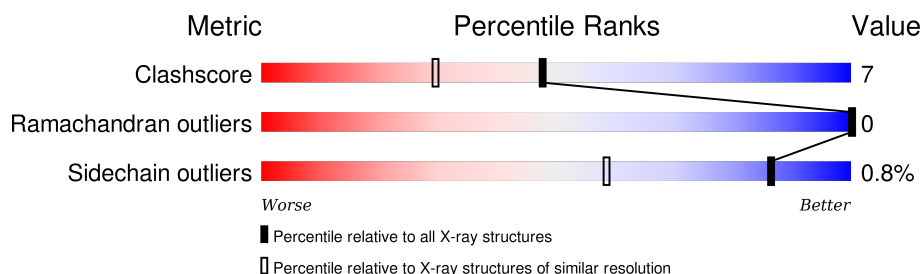
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.10 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	1055 (1.14-1.06)
Ramachandran outliers	100387	1016 (1.14-1.06)
Sidechain outliers	100360	1014 (1.14-1.06)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	76	<div> <div>92%</div> <div>8%</div> </div>
2	B	67	<div> <div>93%</div> <div>7%</div> </div>
3	C	177	<div> <div>90%</div> <div>7%</div> <div>•</div> </div>
3	D	177	<div> <div>93%</div> <div>5%</div> <div>•</div> </div>

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 5197 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Phycoerythrin alpha-3 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	76	Total	C	N	O	S	4	6	0
			592	366	99	123	4			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	4	LYZ	LYS	MODIFIED RESIDUE	UNP Q00433

- Molecule 2 is a protein called Phycoerythrin alpha-2 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	67	Total	C	N	O	S	0	6	0
			520	326	88	100	6			

- Molecule 3 is a protein called B-phycoerythrin beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	173	Total	C	N	O	S	12	16	0
			1330	822	227	271	10			
3	D	177	Total	C	N	O	S	7	12	0
			1344	832	230	271	11			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	50	CYS	VAL	CONFLICT	UNP P27198
C	56	VAL	TYR	CONFLICT	UNP P27198
C	61	CYS	GLU	CONFLICT	UNP P27198
C	65	SER	HIS	CONFLICT	UNP P27198
C	72	MEN	ASN	MODIFIED RESIDUE	UNP P27198
C	73	CYS	GLU	CONFLICT	UNP P27198
D	50	CYS	VAL	CONFLICT	UNP P27198

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Chain	Residue	Modelled	Actual	Comment	Reference
D	56	VAL	TYR	CONFLICT	UNP P27198
D	61	CYS	GLU	CONFLICT	UNP P27198
D	65	SER	HIS	CONFLICT	UNP P27198
D	72	MEN	ASN	MODIFIED RESIDUE	UNP P27198
D	73	CYS	GLU	CONFLICT	UNP P27198

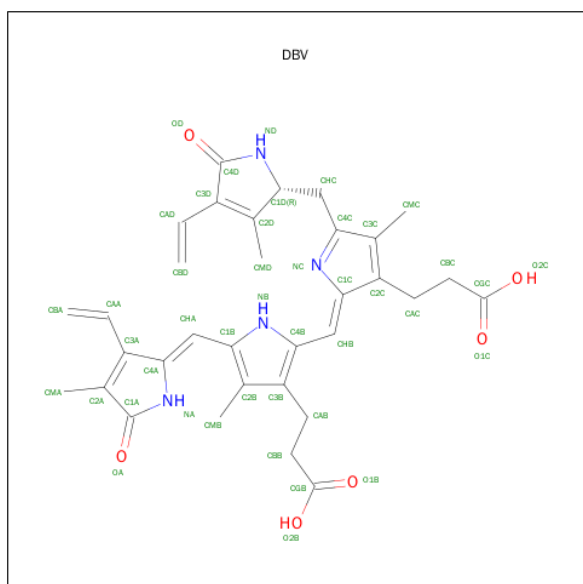
- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	1	Total Cl 1 1	0	0

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Mg 1 1	0	0
5	D	1	Total Mg 1 1	0	0

- Molecule 6 is 15,16-DIHYDROBILIVERDIN (three-letter code: DBV) (formula: C<sub>33</sub>H<sub>36</sub>N<sub>4</sub>O<sub>6</sub>).



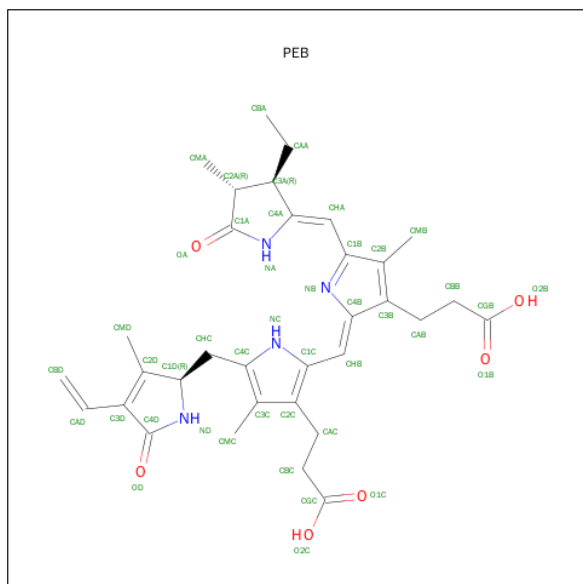
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C N O 43 33 4 6	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	N	O	0	0
			43	33	4	6		

- Molecule 7 is PHYCOERYTHROBILIN (three-letter code: PEB) (formula:  $C_{33}H_{40}N_4O_6$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	C	1	Total	C	N	O	0	0
			43	33	4	6		
7	C	1	Total	C	N	O	0	0
			43	33	4	6		
7	C	1	Total	C	N	O	0	0
			43	33	4	6		
7	D	1	Total	C	N	O	0	0
			43	33	4	6		
7	D	1	Total	C	N	O	0	0
			43	33	4	6		
7	D	1	Total	C	N	O	0	0
			43	33	4	6		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	207	Total O 207 207	0	0
8	B	170	Total O 170 170	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	C	299	Total 299	O 299	0	0
8	D	388	Total 388	O 388	0	0

### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

- Molecule 1: Phycoerythrin alpha-3 chain

Chain A:  92% 8%




- Molecule 2: Phycoerythrin alpha-2 chain

Chain B:  93% 7%



- Molecule 3: B-phycoerythrin beta chain

Chain C:  90% 7% .



- Molecule 3: B-phycoerythrin beta chain

Chain D:  93% 5% .



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	62.98 Å 82.56 Å 89.53 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.86 – 1.10	Depositor
% Data completeness (in resolution range)	98.8 (60.86-1.10)	Depositor
$R_{merge}$	0.03	Depositor
$R_{sym}$	0.03	Depositor
Refinement program	REFMAC 5.2.0003	Depositor
R, $R_{free}$	0.096 , 0.116	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5197	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	10.0	wwPDB-VP



## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, CL, MEN, LYZ, PEB, DBV

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.88	1/610 (0.2%)	0.91	2/812 (0.2%)
2	B	0.83	1/542 (0.2%)	0.85	0/719
3	C	0.82	0/1385	0.89	1/1862 (0.1%)
3	D	0.90	5/1383 (0.4%)	0.93	7/1861 (0.4%)
All	All	0.86	7/3920 (0.2%)	0.90	10/5254 (0.2%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	108	ARG	CG-CD	-6.73	1.35	1.51
2	B	63	ALA	CA-CB	-6.43	1.39	1.52
1	A	74	SER	CA-C	5.83	1.68	1.52
3	D	7	ARG	CD-NE	-5.83	1.36	1.46
3	D	31	ILE	CA-C	5.79	1.68	1.52
3	D	36[A]	LYS	CD-CE	-5.04	1.38	1.51
3	D	36[B]	LYS	CD-CE	-5.04	1.38	1.51

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	77	ARG	NE-CZ-NH2	-9.97	115.31	120.30
3	D	77	ARG	NE-CZ-NH1	8.83	124.71	120.30
3	D	1	MET	CA-CB-CG	6.36	124.11	113.30
3	D	7	ARG	CD-NE-CZ	5.93	131.90	123.60
1	A	13[A]	ILE	CA-CB-CG1	5.91	122.24	111.00
1	A	13[B]	ILE	CA-CB-CG1	5.91	122.24	111.00
3	D	7	ARG	CG-CD-NE	5.65	123.66	111.80
3	D	91	ARG	NE-CZ-NH2	-5.39	117.61	120.30
3	C	78	ARG	NE-CZ-NH1	5.19	122.90	120.30
3	D	91	ARG	NE-CZ-NH1	5.13	122.86	120.30

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	592	0	602	5	1
2	B	520	0	568	4	0
3	C	1330	0	1363	19	0
3	D	1344	0	1380	21	0
4	C	1	0	0	1	0
5	A	1	0	0	0	0
5	D	1	0	0	0	0
6	A	43	0	33	2	0
6	B	43	0	33	2	0
7	C	129	0	110	6	0
7	D	129	0	110	5	0
8	A	207	0	0	7	0
8	B	170	0	0	3	1
8	C	299	0	0	16	1
8	D	388	0	0	13	1
All	All	5197	0	4199	61	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (61) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:149[A]:LYS:HD2	8:D:2113:HOH:O	1.25	1.30
3:D:149[A]:LYS:CD	8:D:2113:HOH:O	1.76	1.27
2:B:34[B]:LYS:HD3	8:B:1194:HOH:O	1.42	1.19
1:A:24[B]:LYS:CE	8:A:2006:HOH:O	1.91	1.17
3:D:149[A]:LYS:CE	8:D:2113:HOH:O	1.91	1.14
8:A:2073:HOH:O	3:D:149[B]:LYS:HE3	1.45	1.14
1:A:24[B]:LYS:HE3	8:A:2006:HOH:O	1.47	1.07
8:A:2073:HOH:O	3:D:149[B]:LYS:CE	2.03	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:149[B]:LYS:NZ	8:C:3107:HOH:O	1.94	0.99
3:D:149[A]:LYS:HE3	8:D:2113:HOH:O	1.59	0.95
3:C:108:ARG:HD2	8:C:3227:HOH:O	1.64	0.95
3:C:111[A]:ASN:OD1	8:C:3291:HOH:O	1.85	0.94
3:C:108:ARG:CD	8:C:3227:HOH:O	2.18	0.88
3:C:149[B]:LYS:CE	8:C:3107:HOH:O	2.21	0.87
2:B:51[A]:THR:HG21	8:B:1159:HOH:O	1.78	0.84
1:A:76:ASN:HB2	8:C:3300:HOH:O	1.78	0.84
2:B:4[A]:LYS:NZ	8:B:1065:HOH:O	2.08	0.83
3:C:115[B]:GLU:CD	8:C:3298:HOH:O	2.17	0.82
1:A:24[B]:LYS:HE2	8:A:2006:HOH:O	1.62	0.79
3:D:36[B]:LYS:HE2	7:D:258:PEB:HMB3	1.71	0.71
3:D:10:THR:HG22	3:D:108:ARG:NH2	2.06	0.69
3:C:115[B]:GLU:OE1	8:C:3163:HOH:O	0.70	0.69
4:C:3001:CL:CL	8:D:2234:HOH:O	2.48	0.67
3:C:108:ARG:NE	8:C:3227:HOH:O	2.26	0.65
8:A:2073:HOH:O	3:D:149[B]:LYS:HE2	1.85	0.63
1:A:9:PRO:O	3:C:91[B]:ARG:HD3	2.00	0.61
6:B:219:DBV:HMB3	6:B:219:DBV:HNA	1.65	0.61
3:D:36[B]:LYS:HE3	8:D:2251:HOH:O	2.01	0.61
3:C:64:PRO:HG3	8:C:3300:HOH:O	2.01	0.60
3:C:176:ILE:HG13	8:C:3249:HOH:O	2.01	0.59
6:A:219:DBV:HMB3	6:A:219:DBV:HNA	1.67	0.59
7:D:282:PEB:HMB2	7:D:282:PEB:HNA	1.68	0.59
7:C:282:PEB:HMB2	7:C:282:PEB:HNA	1.70	0.57
3:C:64:PRO:CG	8:C:3300:HOH:O	2.52	0.56
3:C:91[A]:ARG:HG3	8:C:3219:HOH:O	2.06	0.55
3:D:144:ASN:C	3:D:144:ASN:HD22	2.09	0.54
3:D:115[A]:GLU:CG	8:D:2298:HOH:O	2.56	0.53
3:C:149[B]:LYS:HE3	8:C:3107:HOH:O	2.01	0.52
6:B:219:DBV:HMB3	6:B:219:DBV:NA	2.25	0.52
6:A:219:DBV:HMB3	6:A:219:DBV:NA	2.25	0.51
3:D:115[A]:GLU:HG2	8:D:2298:HOH:O	2.10	0.50
7:D:258:PEB:HNA	7:D:258:PEB:HMB2	1.76	0.50
3:D:36[B]:LYS:CE	7:D:258:PEB:HMB3	2.39	0.50
7:C:258:PEB:HHA1	7:C:258:PEB:HBA3	1.93	0.50
3:D:36[B]:LYS:CE	8:D:2251:HOH:O	2.60	0.49
3:C:149[A]:LYS:HE3	8:C:3107:HOH:O	2.12	0.49
7:C:258:PEB:HNA	7:C:258:PEB:HMB2	1.80	0.47
3:C:31:ILE:HD12	3:C:37:ARG:HD2	1.96	0.47
3:C:149[A]:LYS:NZ	8:C:3109:HOH:O	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:149[B]:LYS:HE2	8:D:2085:HOH:O	2.14	0.46
7:C:250:PEB:HNA	7:C:250:PEB:HMB3	1.80	0.46
3:D:115[A]:GLU:HG3	8:D:2298:HOH:O	2.17	0.45
3:D:149[B]:LYS:CE	8:D:2085:HOH:O	2.65	0.45
3:D:144:ASN:ND2	3:D:146:ALA:H	2.15	0.44
7:C:258:PEB:NA	7:C:258:PEB:HMB2	2.35	0.42
8:A:2019:HOH:O	2:B:66:THR:HB	2.18	0.42
3:D:31:ILE:HD12	3:D:37:ARG:HD2	2.01	0.41
3:D:149[B]:LYS:NZ	8:D:2085:HOH:O	2.45	0.41
3:C:72:MEN:HE22	7:C:282:PEB:HBB2	2.03	0.41
7:D:258:PEB:HHA1	7:D:258:PEB:HBA3	2.03	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37[B]:GLU:OE2	8:B:1187:HOH:O[3_545]	2.17	0.03
8:C:3287:HOH:O	8:D:2254:HOH:O[4_556]	2.19	0.01

## 5.3 Torsion angles

### 5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	79/76 (104%)	77 (98%)	2 (2%)	0	100	100
2	B	71/67 (106%)	69 (97%)	2 (3%)	0	100	100
3	C	185/177 (104%)	184 (100%)	1 (0%)	0	100	100
3	D	186/177 (105%)	185 (100%)	1 (0%)	0	100	100
All	All	521/497 (105%)	515 (99%)	6 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	66/60 (110%)	66 (100%)	0	100	100
2	B	57/51 (112%)	57 (100%)	0	100	100
3	C	154/141 (109%)	152 (99%)	2 (1%)	76	40
3	D	153/141 (108%)	150 (98%)	3 (2%)	63	22
All	All	430/393 (109%)	425 (99%)	5 (1%)	86	43

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	155[A]	GLN
3	C	155[B]	GLN
3	D	144	ASN
3	D	149[A]	LYS
3	D	149[B]	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
3	D	144	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

3 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length

(or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
1	LYZ	A	4	1	8,9,10	1.18	1 (12%)	6,10,12	0.99	1 (16%)
3	MEN	C	72	3	7,8,9	0.81	0	5,9,11	1.37	1 (20%)
3	MEN	D	72	3	7,8,9	0.78	0	5,9,11	1.34	1 (20%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	LYZ	A	4	1	-	0/7/9/11	0/0/0/0
3	MEN	C	72	3	-	0/6/8/10	0/0/0/0
3	MEN	D	72	3	-	0/6/8/10	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	4	LYZ	CB-CA	-2.88	1.51	1.53

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	72	MEN	O-C-CA	-2.94	117.82	125.49
3	C	72	MEN	O-C-CA	-2.61	118.70	125.49
1	A	4	LYZ	O-C-CA	-2.18	119.80	125.49

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	72	MEN	1	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 3 are monoatomic - leaving 8 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
6	DBV	A	219	1	32,46,46	1.53	2 (6%)	34,67,67	0.99	2 (5%)
6	DBV	B	219	2	32,46,46	1.99	4 (12%)	34,67,67	1.84	2 (5%)
7	PEB	C	250	3	36,46,46	1.92	5 (13%)	38,67,67	1.24	3 (7%)
7	PEB	C	258	3	36,46,46	1.70	4 (11%)	38,67,67	1.43	7 (18%)
7	PEB	C	282	3	36,46,46	1.34	5 (13%)	38,67,67	1.58	6 (15%)
7	PEB	D	250	3	36,46,46	1.78	6 (16%)	38,67,67	1.26	5 (13%)
7	PEB	D	258	3	36,46,46	1.44	6 (16%)	38,67,67	1.54	4 (10%)
7	PEB	D	282	3	36,46,46	1.48	5 (13%)	38,67,67	1.61	5 (13%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	DBV	A	219	1	-	0/21/74/74	0/4/4/4
6	DBV	B	219	2	-	0/21/74/74	0/4/4/4
7	PEB	C	250	3	-	2/19/74/74	0/4/4/4
7	PEB	C	258	3	-	2/19/74/74	0/4/4/4
7	PEB	C	282	3	-	2/19/74/74	0/4/4/4
7	PEB	D	250	3	-	2/19/74/74	0/4/4/4
7	PEB	D	258	3	-	2/19/74/74	0/4/4/4
7	PEB	D	282	3	-	2/19/74/74	0/4/4/4

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	C	258	PEB	C2A-C1A	-6.11	1.46	1.52
7	D	258	PEB	CAA-C3A	-3.39	1.47	1.54
7	C	258	PEB	C1A-NA	-2.58	1.34	1.37
7	D	282	PEB	C2A-C1A	-2.53	1.49	1.52
7	D	250	PEB	C1B-C2B	-2.41	1.40	1.45
6	B	219	DBV	CMA-C2A	-2.25	1.46	1.50
7	C	282	PEB	C4B-C3B	-2.17	1.41	1.45
7	C	250	PEB	C4B-C3B	-2.15	1.41	1.45
7	D	282	PEB	C4B-C3B	-2.14	1.41	1.45
7	D	250	PEB	C4B-C3B	-2.08	1.42	1.45
7	D	258	PEB	CBB-CAB	-2.06	1.42	1.52
7	D	258	PEB	C1B-C2B	-2.04	1.41	1.45
7	C	282	PEB	C2A-C1A	-2.02	1.50	1.52
7	D	282	PEB	C3B-C2B	2.07	1.41	1.36
7	D	258	PEB	C3B-C2B	2.08	1.41	1.36
6	B	219	DBV	CHA-C4A	2.11	1.39	1.34
7	D	258	PEB	CHA-C1B	2.19	1.45	1.40
7	D	250	PEB	CHA-C1B	2.20	1.45	1.40
7	C	250	PEB	C3B-C2B	2.24	1.41	1.36
7	C	282	PEB	C3B-C2B	2.25	1.41	1.36
7	C	282	PEB	CHA-C1B	2.27	1.46	1.40
7	C	258	PEB	C3B-C2B	2.36	1.41	1.36
7	D	250	PEB	C3B-C2B	2.38	1.41	1.36
7	C	250	PEB	CHA-C1B	2.61	1.46	1.40
7	D	282	PEB	CHA-C1B	2.68	1.47	1.40
7	C	250	PEB	CBD-CAD	4.46	1.52	1.30
7	C	282	PEB	CHB-C4B	4.60	1.39	1.35
7	D	250	PEB	CBD-CAD	4.83	1.54	1.30
6	B	219	DBV	CBA-CAA	4.89	1.54	1.30
6	A	219	DBV	CBA-CAA	4.96	1.54	1.30
6	A	219	DBV	CHB-C1C	5.29	1.39	1.35
7	C	258	PEB	CHB-C4B	5.54	1.40	1.35
7	D	258	PEB	CHB-C4B	5.69	1.40	1.35
7	D	282	PEB	CHB-C4B	6.31	1.40	1.35
7	D	250	PEB	CHB-C4B	7.03	1.41	1.35
6	B	219	DBV	CHB-C1C	8.24	1.42	1.35
7	C	250	PEB	CHB-C4B	9.05	1.43	1.35

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	282	PEB	CAC-CBC-CGC	-3.74	105.89	112.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	282	PEB	CHA-C1B-NB	-3.73	117.80	124.91
7	D	282	PEB	CHA-C1B-NB	-3.71	117.84	124.91
7	D	258	PEB	CHA-C1B-NB	-3.57	118.10	124.91
7	C	258	PEB	OA-C1A-NA	-3.22	120.93	124.83
7	D	250	PEB	CHA-C1B-NB	-3.18	118.86	124.91
7	D	250	PEB	CBD-CAD-C3D	-3.03	111.64	127.01
7	C	250	PEB	CBD-CAD-C3D	-2.77	112.98	127.01
6	B	219	DBV	CBA-CAA-C3A	-2.68	113.42	127.01
7	D	250	PEB	CAB-CBB-CGB	-2.59	108.00	112.75
7	D	258	PEB	CAC-CBC-CGC	-2.51	108.15	112.75
6	A	219	DBV	CBA-CAA-C3A	-2.49	114.39	127.01
7	C	258	PEB	CHA-C1B-NB	-2.42	120.30	124.91
7	C	282	PEB	CAB-C3B-C2B	-2.25	124.08	128.01
7	C	258	PEB	CMC-C3C-C4C	-2.24	124.88	127.14
7	C	250	PEB	CHA-C1B-NB	-2.14	120.84	124.91
7	D	282	PEB	C3B-C4B-NB	-2.10	106.59	109.86
6	A	219	DBV	C2C-C1C-NC	-2.06	106.67	109.86
7	C	258	PEB	C4B-NB-C1B	2.02	110.51	106.51
7	D	250	PEB	C2A-C1A-NA	2.04	110.25	108.30
7	C	258	PEB	CAA-C3A-C4A	2.20	118.32	112.67
7	C	258	PEB	CHA-C1B-C2B	2.59	131.19	124.88
7	C	282	PEB	CMB-C2B-C1B	2.59	129.28	125.06
7	D	250	PEB	CHA-C1B-C2B	2.69	131.43	124.88
7	D	282	PEB	CMB-C2B-C1B	2.83	129.66	125.06
7	C	282	PEB	CHA-C1B-C2B	3.11	132.46	124.88
7	D	258	PEB	CHA-C1B-C2B	3.11	132.47	124.88
7	C	250	PEB	CAB-CBB-CGB	3.14	118.50	112.75
7	D	282	PEB	CHA-C1B-C2B	3.33	133.02	124.88
7	C	282	PEB	C2A-C1A-NA	3.35	111.50	108.30
7	C	258	PEB	C2A-C1A-NA	4.15	112.26	108.30
7	D	282	PEB	CAB-CBB-CGB	4.74	121.42	112.75
7	D	258	PEB	C3A-C2A-C1A	5.65	108.15	103.41
6	B	219	DBV	CAC-CBC-CGC	9.17	129.56	112.75

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	D	250	PEB	C4A-CHA-C1B-C2B
7	C	250	PEB	C4A-CHA-C1B-C2B
7	C	258	PEB	C4A-CHA-C1B-C2B
7	D	258	PEB	C4A-CHA-C1B-C2B

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Mol	Chain	Res	Type	Atoms
7	C	282	PEB	C4A-CHA-C1B-C2B
7	C	250	PEB	C4A-CHA-C1B-NB
7	D	250	PEB	C4A-CHA-C1B-NB
7	D	282	PEB	C4A-CHA-C1B-C2B
7	C	258	PEB	C4A-CHA-C1B-NB
7	D	258	PEB	C4A-CHA-C1B-NB
7	D	282	PEB	C4A-CHA-C1B-NB
7	C	282	PEB	C4A-CHA-C1B-NB

There are no ring outliers.

7 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	219	DBV	2	0
6	B	219	DBV	2	0
7	C	250	PEB	1	0
7	C	258	PEB	3	0
7	C	282	PEB	2	0
7	D	258	PEB	4	0
7	D	282	PEB	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.